

**SUMMARY OF ANALYSIS (SAMPLE ID: SA35642)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Order ID:</b> OR10532 <b>Lot Number:</b> M00065C13246697663 <b>Production Run:</b> E20230714CBSSLR01	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4grams <b>Date Collected:</b> 07/17/2023 <b>Date Received:</b> 07/17/2023 <b>Date Completed:</b> 07/20/2023
<b>Cultivar (Strain) or Sample Description:</b> Crunch Berries Solventless Live Hash Rosin 1g Jar			

\*This page is simply a summary of the analysis performed. For analytical details, please consult the individual Certificate(s) of Analysis for each of the specific test(s) performed. All contaminant action levels are referenced from the State of Arkansas MMJ testing guidelines.

**Moisture Content (%)**

Not Tested

**Water Activity (aw)**

Not Tested

**PASS/FAIL**

**PASS**

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

<b>Cannabinoids (Top 3)</b>	<b>(%)</b>	<b>mg/g</b>
THCA	83.0	830
Δ9-THC	3.39	33.9
CBGA	1.97	19.7
TOTAL CBD	-	-
TOTAL THC	76.2	762
TOTAL CANNABINOIDS	88.9	889

<b>Terpenes (Top 5)</b>	<b>(%)</b>	<b>µg/g</b>
α-Pinene	3.14	31400
β-Pinene	1.83	18300
β-Caryophyllene	0.516	5160
cis-Nerolidol	0.360	3600
d-Limonene	0.263	2630
TOTAL TERPENES	6.59	65900

<b>Contaminants</b>	<b>PASS/FAIL</b>
Heavy Metals:	<b>PASS</b>
Microbiology:	<b>PASS</b>
Pesticides:	<b>PASS</b>
Residual Solvents:	<b>PASS</b>

**Sample Picture Upon Receipt**



Scan the QR code to verify results.

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*Kyle W. Felling*  
Kyle W. Felling, Ph.D.  
Laboratory Director

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**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35642)**

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**CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)**

**Analysis Date/Time:** 07/18/2023 1927  
**Analyst:** PW

**Method:** HPLC/DAD  
**Instrument:** Agilent 1100

**Moisture Content (%):** -  
**Water Activity (aw):** -

<u>Cannabinoid</u>	<u>Result (%)</u>	<u>Result (mg/g)</u>	<u>LOD (mg/g)</u>	<u>LOQ (mg/g)</u>	<u>Result (mg/mL)</u>	<u>Per Serving (mg)</u>	<u>Per Unit (mg)</u>
CBC	ND	ND	0.111	0.259	-	-	-
CBCA	ND	ND	0.345	0.804	-	-	-
CBD	ND	ND	0.783	1.83	-	-	-
CBDa	ND	ND	0.288	0.673	-	-	-
CBDV	ND	ND	0.126	0.293	-	-	-
CBDVA	ND	ND	0.335	0.782	-	-	-
CBG	0.169	1.69	0.508	1.19	-	1.69	1.69
CBGA	1.97	19.7	0.720	0.850	-	19.7	19.7
CBL	ND	ND	0.587	1.37	-	-	-
CBN	ND	ND	0.270	0.630	-	-	-
CBNA	ND	ND	0.291	0.678	-	-	-
Δ9-THC	3.39	33.9	0.323	0.753	-	33.9	33.9
Δ8-THC	ND	ND	0.504	1.18	-	-	-
THCA	83.0	830	0.175	0.410	-	830	830
THCV	ND	ND	0.420	0.980	-	-	-
THCVA	0.427	4.27	0.134	0.312	-	4.27	4.27
<b>TOTAL</b>	<b>88.9</b>	<b>889</b>				<b>889</b>	<b>889</b>
<b>TOTAL CBC</b>	-	-				-	-
<b>TOTAL CBD</b>	-	-				-	-
<b>TOTAL CBDV</b>	-	-				-	-
<b>TOTAL CBG</b>	1.90	19.0				19.0	19.0
<b>TOTAL CBN</b>	-	-				-	-
<b>TOTAL THC</b>	76.2	762				762	762
<b>TOTAL THCv</b>	0.370	3.70				3.70	3.70



**SERVING MASS (g):** 1.00  
**SERVINGS/UNIT:** 1

"-" Not detected above LOD.

*Deviations from standard operating procedure:*  
None

*Recoveries for all analyte standards:* 90-110%  
*Replicate Uncertainties:* <5% RSD, <20% RPD  
*Sample/Reagent Blanks:* <RL for all analytes

Values for plant matter are adjusted for moisture content.

Total CBC = (CBCA x 0.877) + CBC  
Total CBD = (CBDa x 0.877) + CBD  
Total CBDV = (CBDVA x 0.867) + CBDV  
Total CBG = (CBGA x 0.878) + CBG  
Total CBN = (CBNA x 0.876) + CBN  
Total THC = (THCA x 0.877) + Δ9-THC  
Total THCv = (THCVA x 0.867) + THCv

Percentage results are reported by mass.

mg/g results are reported as mass component per mass material.

*Abbreviations:* DAD - Diode Array Detector, HPLC - High Pressure Liquid Chromatography, RL - Reporting Limit, RPD - Relative Percent Difference, RSD - Relative Standard Deviation, DET - Detected (less than LOQ), LOD - Limit of Detection, LOQ - Limit of Quantitation, UM - Measurement Uncertainty

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<b>Cultivar (Strain) or Sample Description:</b> Crunch Berries Solventless Live Hash Rosin 1g Jar			<b>Date Completed:</b> 07/20/2023

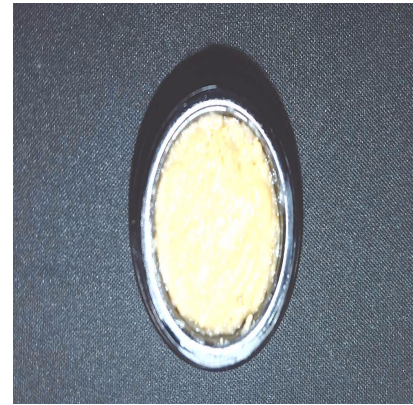
**TERPENOID PROFILE**

**Analysis Date/Time:** 07/19/2023 1708  
**Analyst:** KF

**Method:** GC/MS  
**Instrument:** Agilent 7890/5975

**Deviations from SOP:**  
None

<u>Terpene</u>	<u>Result (µg/g)</u>	<u>Result (%)</u>
α-Bisabolol	ND	-
Camphene	584	0.0584
δ-3-Carene	ND	-
β-Caryophyllene	5160	0.516
Caryophyllene oxide	ND	-
p-Cymene	ND	-
Eucalyptol	ND	-
Geraniol	ND	-
Guaiol	ND	-
α-Humulene	2140	0.214
Isopulegol	ND	-
d-Limonene	2630	0.263
Linalool	ND	-
β-Myrcene	2040	0.204
cis-Nerolidol	3600	0.360
trans-Nerolidol	ND	-
α-Ocimene	ND	-
β-Ocimene	ND	-
α-Pinene	31400	3.14
β-Pinene	18300	1.83
α-Terpinene	ND	-
γ-Terpinene	ND	-
Terpinolene	ND	-
<b>TOTAL</b>	<b>65900</b>	<b>6.59</b>



*Abbreviations:* GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit  
*Abbreviations:* ND - Not Detected, LOD - Limit of Detection, LOQ - Limit of Quantitation

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**Reporting Limit (µg/g): 243**

"-" Not detected above LOD.

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Kyle W. Felling, Ph.D.  
Laboratory Director



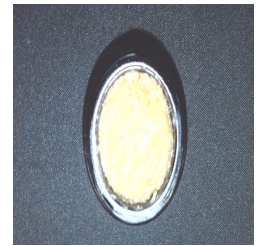
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**RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)**

<b>Analysis Date/Time:</b> 07/19/2023 0320	<b>Method:</b> HS/GC/MS	<b>Deviations from SOP:</b>
<b>Analyst:</b> KF	<b>Instrument:</b> Agilent 7890/5975	None

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	96.3	193	5000	n-Heptane (142-82-5)	-	96.3	193	5000
Acetonitrile (75-5-8)	-	96.3	193	410	n-Hexane (110-54-3)	-	33.7	67.4	290
Benzene (71-43-2)	-	0.963	1.93	2	Isobutane (75-28-5)	-	96.3	193	5000
n-Butane (106-97-2)	-	96.3	193	5000	Isopropanol (67-63-0)	-	96.3	193	5000
1-Butanol (71-36-3)	-	96.3	193	5000	Isopropyl acetate (108-21-4)	-	96.3	193	5000
2-Butanol (78-92-2)	-	96.3	193	5000	Isopropyl benzene (98-82-8)	-	9.63	19.3	70
2-Butanone (78-93-3)	-	96.3	193	5000	Methanol (67-56-1)	-	96.3	193	3000
Cyclohexane (110-82-7)	-	96.3	193	3880	2-Methylbutane (78-78-4)	-	96.3	193	5000
1,2-Dimethoxyethane (110-71-4)	-	9.63	19.3	100	Methylene chloride (75-9-2)	-	96.3	193	600
N,N-Dimethylacetamide (127-19-5)	-	96.3	193	1090	2-Methylpentane (107-83-5)	-	33.7	67.4	290
2,2-Dimethylbutane (75-83-2)	-	33.7	67.4	290	3-Methylpentane (96-10-0)	-	33.7	67.4	290
2,3-Dimethylbutane (79-29-8)	-	33.7	67.4	290	n-Pentane (109-66-0)	-	96.3	193	5000
N,N-Dimethylformamide (68-12-2)	-	96.3	193	880	1-Pentanol (71-41-0)	-	96.3	193	5000
Dimethylsulfoxide (67-68-5)	-	96.3	193	5000	n-Propane (74-98-6)	-	96.3	193	5000
1,4-Dioxane (123-91-1)	-	96.3	193	380	1-Propanol (71-23-8)	-	96.3	193	5000
Ethanol (64-17-5)	-	96.3	193	5000	Pyridine (110-86-1)	-	33.7	67.4	200
2-Ethoxyethanol (110-80-5)	-	33.7	67.4	160	Tetrahydrofuran (109-99-9)	-	96.3	193	720
Ethyl ether (60-29-7)	-	96.3	193	5000	Tetramethylene sulfone (126-33-0)	-	33.7	67.4	160
Ethyl acetate (141-78-6)	-	96.3	193	5000	Toluene (108-88-3)	-	96.3	193	890
Ethyl benzene (100-41-4)	-	96.3	193	2170	o-Xylene (95-47-6)	-	96.3	193	2170
Ethylene glycol (107-21-1)	-	96.3	193	620	m,p-Xylene (108-38-3 or 106-42-3)	-	96.3	193	2170
Ethylene oxide (75-21-8)	-	9.63	19.3	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



**Color Key**

RESULT < AL
RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

"\*" - o,m,p-Xylene and Ethylbenzene

Action levels are referenced from the State of Arkansas MMJ testing guidelines.

A value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above.

Solvent	Synonym(s)	Solvent	Synonym(s)
Acetonitrile	Methyl Cyanide, ACN	Ethylene glycol	1,2-Ethanediol
1-Butanol	n-Butanol, Butyl Alcohol	Isobutane	2-Methylpropane
2-Butanol	sec-Butyl alcohol	Isopropanol	2-Propanol, IPA
2-Butanone	Methyl ethyl ketone, MEK	Isopropyl Acetate	Acetic acid isopropyl ester
1,2-Dimethoxyethane	Monoglyme	Methanol	Methyl alcohol
2,3-Dimethylbutane	Neohexane	2-Methylbutane	Isopentane
2,3-Dimethylbutane	Diisopropyl	Methylene chloride	Dichloromethane
N,N-Dimethylformamide	DMF	2-Methylpentane	Isohexane
Dimethylsulfoxide	DMSO	1-Pentanol	n-Amyl alcohol
2-Ethoxyethanol	Cellosolve, Ethyl glycol	1-Propanol	Propyl alcohol
Ethyl ether	Diethyl ether, Ether	Tetrahydrofuran	THF
Ethyl acetate	EtOAc	Tetramethylene sulfone	Sulfolane
Ethyl benzene	Phenylethane	Xylene	Dimethylbenzene

Abbreviations: HS-Headspace, GC-Gas Chromatography, MS-Mass Spectrometry, RL-Reporting Limit, AL-Action Level, CAS-Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

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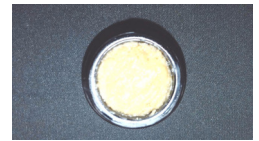
**PESTICIDES PROFILE (SOP: SOP-PEST-001)**

**Analysis Date/Time:** 07/18/2023 2124  
**Analyst:** KF

**Method:** LC/MS/MS  
**Instrument:** Shimadzu LC-8050

**Deviations from SOP:**  
None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0485	0.388	0.5	Kresoxim-methyl (143390-89-0)	-	0.0485	0.388	0.4
Acephate (30560-19-1)	-	0.0485	0.388	0.4	Malathion (121-75-5)	-	0.0485	0.388	0.2
Acequinocyl (57960-19-7)	-	0.0485	0.388	2	Metalaxyl (57837-19-1)	-	0.0485	0.388	0.2
Acetamiprid (135410-20-7)	-	0.0485	0.388	0.2	Methiocarb (2032-65-7)	-	0.0485	0.388	0.2
Aldicarb (116-06-3)	-	0.0485	0.388	0.4	Methomyl (16752-77-5)	-	0.0485	0.388	0.4
Azoxystrobin (131860-33-8)	-	0.0485	0.388	0.2	Methyl parathion (298-0-0)	-	0.0485	0.388	0.2
Bifenazate (149877-41-8)	-	0.0485	0.388	0.2	MGK 264 (113-48-4)	-	0.0485	0.388	0.2
Bifenthrin (82657-04-3)	-	0.0485	0.388	0.2	Myclobutanil (88671-89-0)	-	0.0485	0.388	0.2
Boscalid (188425-85-6)	-	0.0485	0.388	0.4	Naled (300-76-5)	-	0.0485	0.388	0.5
Carbaryl (63-25-2)	-	0.0485	0.388	0.2	Oxamyl (23135-22-0)	-	0.0485	0.388	1
Carbofuran (1563-66-2)	-	0.0485	0.388	0.2	Paclotrazol (76738-62-0)	-	0.0485	0.388	0.4
Chlorantraniliprole (800008-45-7)	-	0.0485	0.388	0.2	Permethrins (52645-53-1)	-	0.0485	0.388	0.2
Chlorfenapyr (122453-73-0)	-	0.0485	0.388	1	Phosmet (732-11-6)	-	0.0485	0.388	0.2
Chlorpyrifos (2921-88-2)	-	0.0485	0.388	0.2	Piperonyl butoxide (51-03-6)	-	0.0485	0.388	2
Clofentezine (74115-24-5)	-	0.0485	0.388	0.2	Prallethrins (2331-36-9)	-	0.0485	0.388	0.2
Cyfluthrin (68359-37-5)	-	0.0485	0.388	1	Propiconazole (60207-90-1)	-	0.0485	0.388	0.4
Cypermethrin (52315-07-8)	-	0.0485	0.388	1	Propoxur (114-26-1)	-	0.0485	0.388	0.2
Daminozide (1596-84-5)	-	0.0485	0.388	1	Pyrethrins (8003-34-7)	-	0.0485	0.388	1
DDVP (62-73-7)	-	0.0485	0.388	0.1	Pyridaben (96489-71-3)	-	0.0485	0.388	0.2
Diazinon (333-41-5)	-	0.0485	0.388	0.2	Spinosad (168316-95-8)	-	0.0485	0.388	0.2
Dimethoate (60-51-5)	-	0.0485	0.388	0.2	Spiromesifen (283594-90-1)	-	0.0485	0.388	0.2
Ethoprophos (13194-48-4)	-	0.0485	0.388	0.2	Spirotetramat (203313-25-1)	-	0.0485	0.388	0.2
Etofenprox (80844-07-1)	-	0.0485	0.388	0.4	Spiroxamine (118134-30-8)	-	0.0485	0.388	0.4
Etoxazole (153233-91-1)	-	0.0485	0.388	0.2	Tebuconazole (80443-41-0)	-	0.0485	0.388	0.4
Fenoxycarb (72490-01-8)	-	0.0485	0.388	0.2	Thiacloprid (111988-49-9)	-	0.0485	0.388	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0485	0.388	0.4	Thiamethoxam (153719-23-4)	-	0.0485	0.388	0.2
Fipronil (120068-37-3)	-	0.0485	0.388	0.4	Trifloxystrobin (141517-21-7)	-	0.0485	0.388	0.2
Fonicamid (158062-67-0)	-	0.0485	0.388	1					
Fludioxinil (131341-86-1)	-	0.0485	0.388	0.4					
Hexythiazox (78587-05-0)	-	0.0485	0.388	1					
Imazalil (35554-44-0)	-	0.0485	0.388	0.2					
Imidacloprid (138261-41-3)	-	0.0485	0.388	0.4					



**Color Key**

RESULT < AL
RESULT > AL

"DET" detected less than LOQ  
"-." not detected above LOD  
Permethrins measured as the cumulative residue of the *cis*- and *trans*- permethrin isomers.  
Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.  
Action levels are referenced from the State of Arkansas MMJ testing guidelines.  
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Abbreviations: LC - Liquid Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, CAS - Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

Pesticide	Synonym(s)	Pesticide	Synonym(s)	Pesticide	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Sythane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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*Kyle W. Felling*  
Kyle W. Felling, Ph.D.  
Laboratory Director

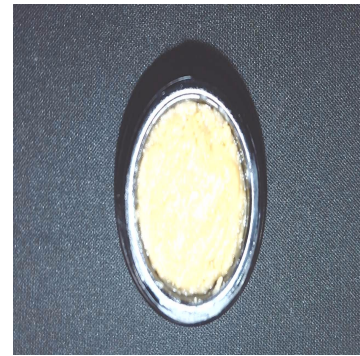
**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35642)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Order ID:</b> OR10532 <b>Lot Number:</b> M00065C13246697663 <b>Production Run:</b> E20230714CBSSLR01	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4grams <b>Date Collected:</b> 07/17/2023 <b>Date Received:</b> 07/17/2023
<b>Cultivar (Strain) or Sample Description:</b> Crunch Berries Solventless Live Hash Rosin 1g Jar			<b>Date Completed:</b> 07/20/2023

**HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)**

<b>Analysis Date/Time:</b> 07/19/2023 1631 (ICP/OES)	<b>Method:</b> ICP/MS	<b>Deviations from SOP:</b>
<b>Analysis Date/Time:</b> - (DMA)	<b>Instrument:</b> Agilent 7500ce	None
<b>Analyst:</b> KF		

<u>Heavy Metal</u>	<u>Result (µg/kg)</u>	<u>LOD (µg/kg)</u>	<u>LOQ (µg/kg)</u>	<u>Action Level (µg/kg)</u>
Arsenic (As)	-	57.4	90.8	200
Cadmium (Cd)	-	57.4	90.8	200
Lead (Pb)	-	57.4	90.8	500
Mercury (Hg)	-	57.4	90.8	100



Abbreviations: ICP - Inductively-Coupled Plasma, OES - Optical Emission Spectroscopy, DMA - Direct Mercury Analyzer, RL - Reporting Limit, AL - Action Level, LOD - Limit of Detection, LOQ - Limit of Quantitation

**Color Key**

RESULT < AL
RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

**Disclaimer:** This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, or prevent any medical-related symptoms. The statements and results herein have not been approved and/or endorsed by the FDA.

**REPORT OF LABORATORY ANALYSIS**

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**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35642)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	<b>Customer ID:</b> 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	<b>Sample ID:</b> SA35642 <b>Lot Number:</b> M00065C13246697663 <b>Production Run:</b> E20230714CBSSLR01	<b>Sample Type:</b> Primary <b>Matrix:</b> Concentrate <b>Mass:</b> 4grams <b>Date Collected:</b> 07/17/2023 <b>Date Received:</b> 07/17/2023
<b>Cultivar (Strain) or Sample Description:</b> Crunch Berries Solventless Live Hash Rosin 1g Jar			<b>Date Completed:</b> 07/20/2023

**MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)**

**Analysis Date/Time:** 07/19/2023 1237      **Method:** Hardy Diagnostics CompactDry      **Deviations from SOP:**  
**Analyst:** PW      **Instrument:** Thermo Incubator      None

<b>Bacteria/Microbe</b>	<b>Result (CFU/g)</b>	<b>Action Level (CFU/g)</b>
Aerobic Plate Count	NT	-
Coliforms, Total	Absent	1
Escherichia Coli (E. Coli)	Absent	100
Mold/Yeast	NT	-
Pseudomonas aeruginosa	NT	-
Salmonella spp.	NT	-
Staphylococcus aureus	NT	-



**Abbreviations:** EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit,  
 AL - Action Level, TNTC - Too Numerous To Count, NT - Not Tested  
 Absent - Not Detected Above RL, Present - Detected Above RL

**Color Key**

**RESULT < AL**

**RESULT > AL**

**Reporting Limit (CFU/g)**  
1

Action levels for microbiology are referenced from the State of Arkansas MMJ testing guidelines.  
 A value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above.

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